

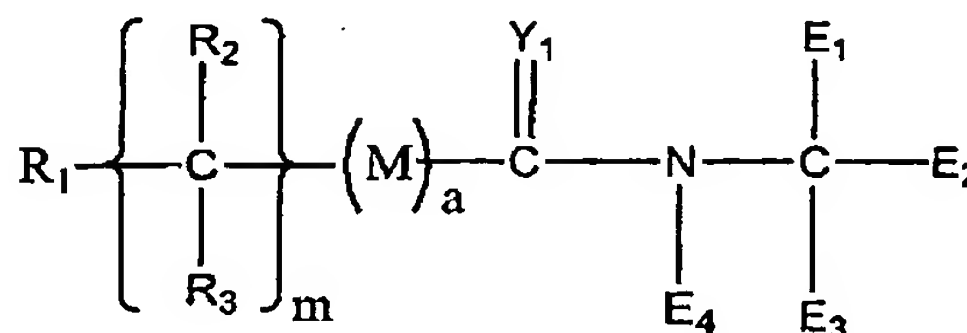
A. AMENDMENTS TO THE CLAIMS:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently amended) A compound comprising the formula:

(I)



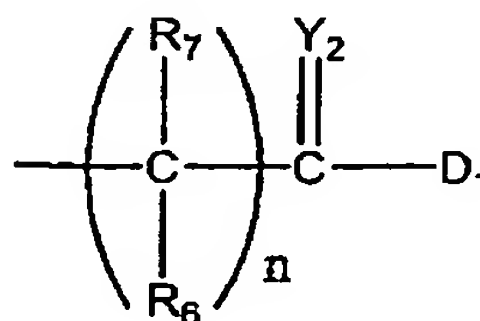
wherein:

R_1 is a polymeric residue;

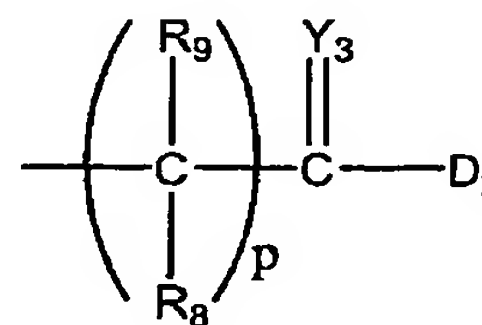
Y_1 is O, S or NR_4 ;

M is O, S or NR_5 ;

E_1 is



E_{2-4} are independently H, E_1 or



(a) is zero or one;

(m) is zero or a positive integer;

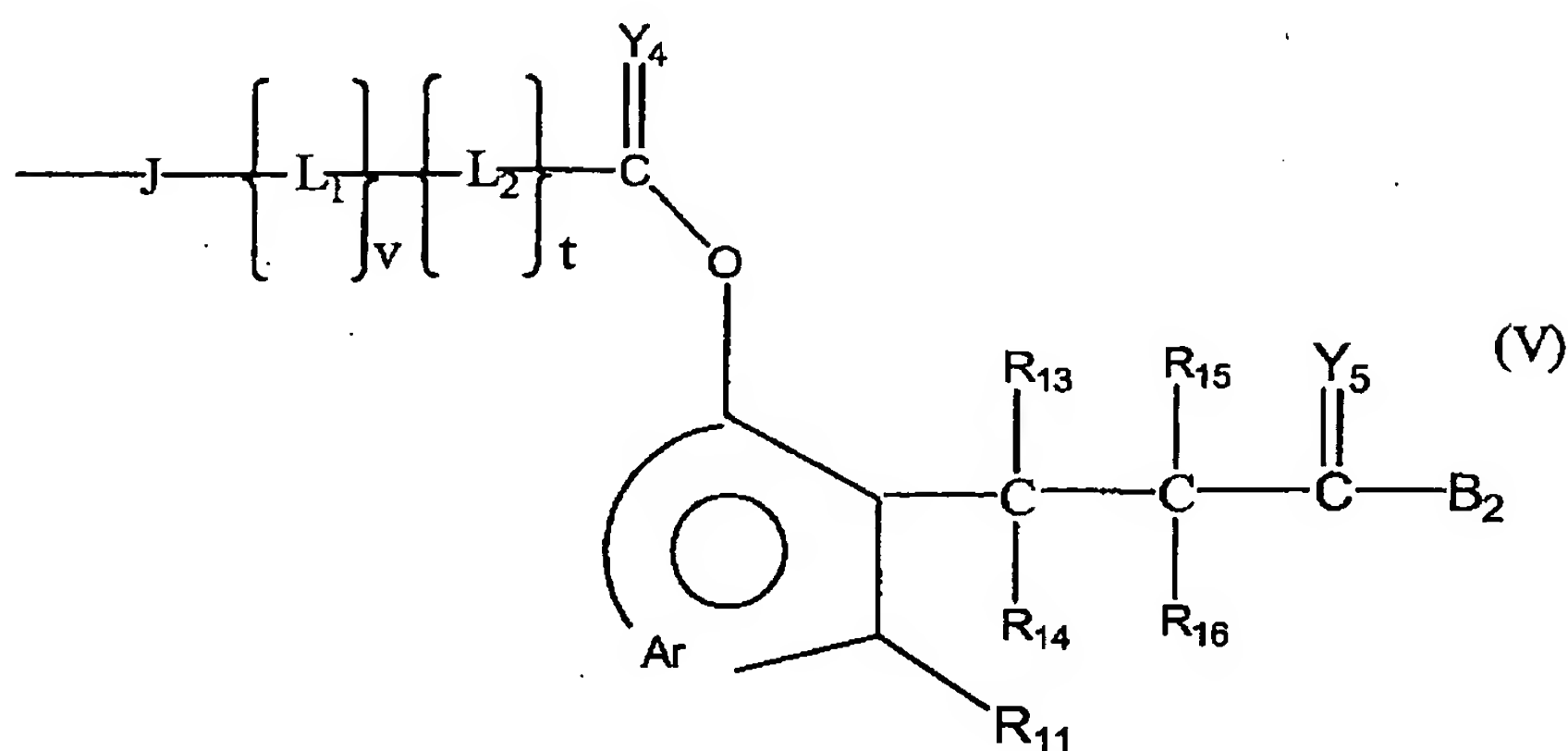
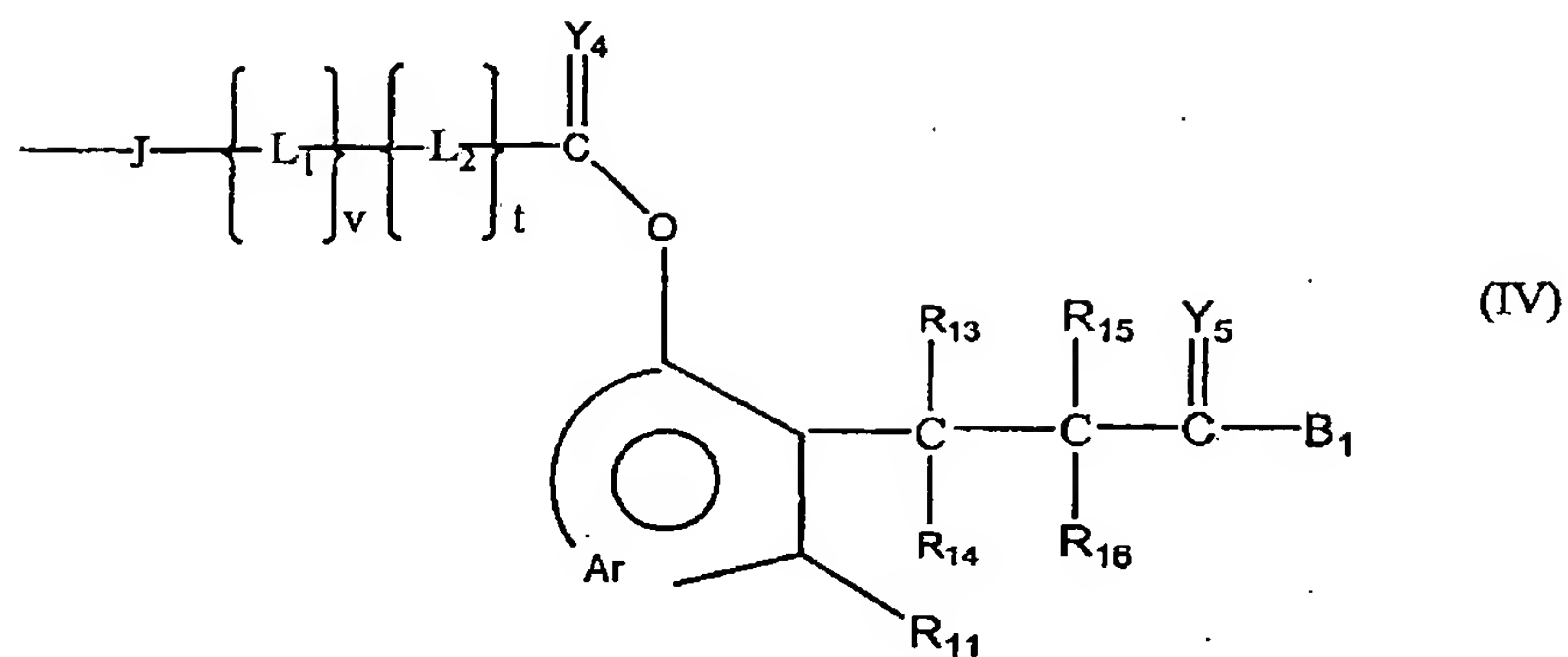
(n) and (p) are independently 0 or a positive integer;

$Y_{2,3}$ are independently O, S or NR_{10} ;

$R_{2,10}$ are independently selected from the group consisting of hydrogen,

C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} hetero-alkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

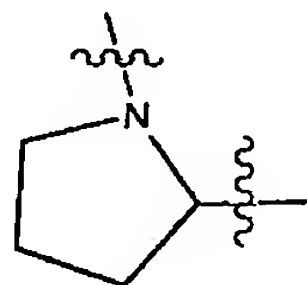
D_1 and D_2 are independently OH,



or a terminal branching group;

wherein (v) and (t) are independently 0 or a positive integer up to about 6;

J is NR_{12} or



L_1 and L_2 are independently selected bifunctional linkers;

$\text{Y}_{4,7}$ are independently selected from the group consisting of O, S and NR_{14} ;

R_{11-14} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls,

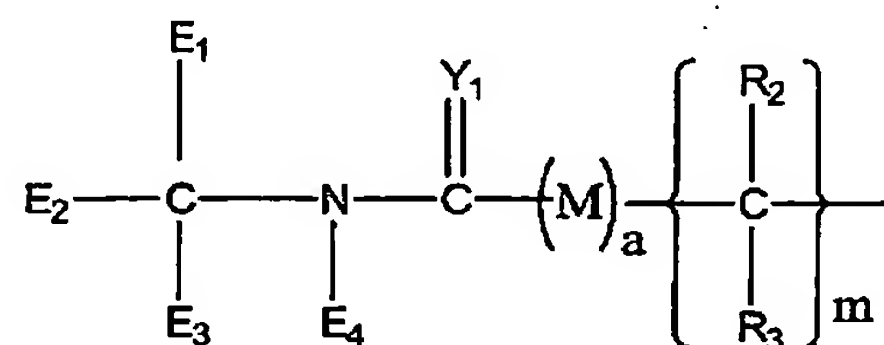
C₃₋₁₂ branched alkyls, C₃₋₈ cycloalkyls, C₁₋₆ substituted alkyls, C₃₋₈ substituted cycloalkyls, aryls, substituted aryls, aralkyls, C₁₋₆ heteroalkyls, substituted C₁₋₆ heteroalkyls, C₁₋₆ alkoxy, phenoxy and C₁₋₆ heteroalkoxy;

Ar is a moiety which when included in Formula (I) forms a multi-substituted aromatic hydrocarbon or a multi-substituted heterocyclic group;

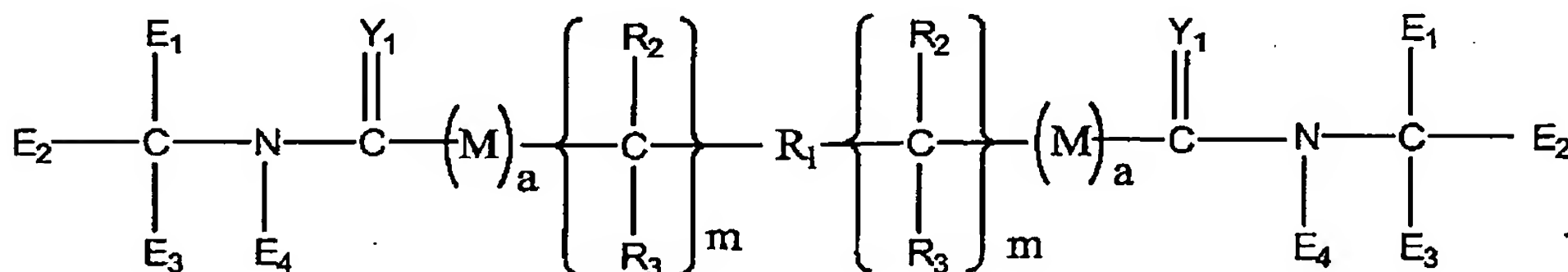
B₁ and B₂ are independently selected from the group consisting of leaving groups, OH, residues of hydroxyl-containing moieties or amine-containing moieties;

provided that E₂₋₄ are not all H.

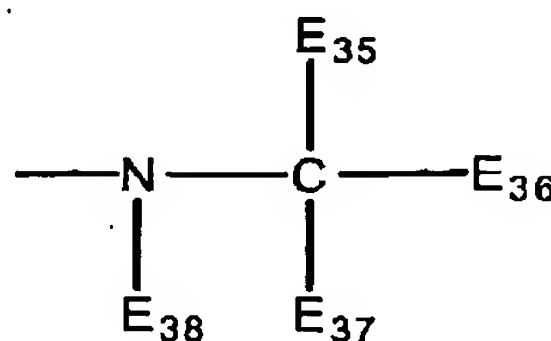
2. (Original) The compound of claim 1, wherein R₁ further comprises a capping group A, selected from the group consisting of hydrogen, NH₂, OH, CO₂H, C₁₋₆ moieties and



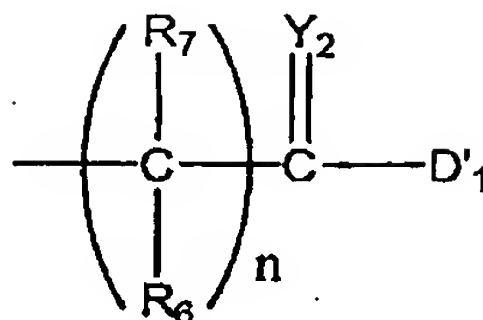
3. (Original) A compound of claim 2, comprising the formula:



4. (Original) The compound of claim 1, wherein said terminal branching group comprises the formula:

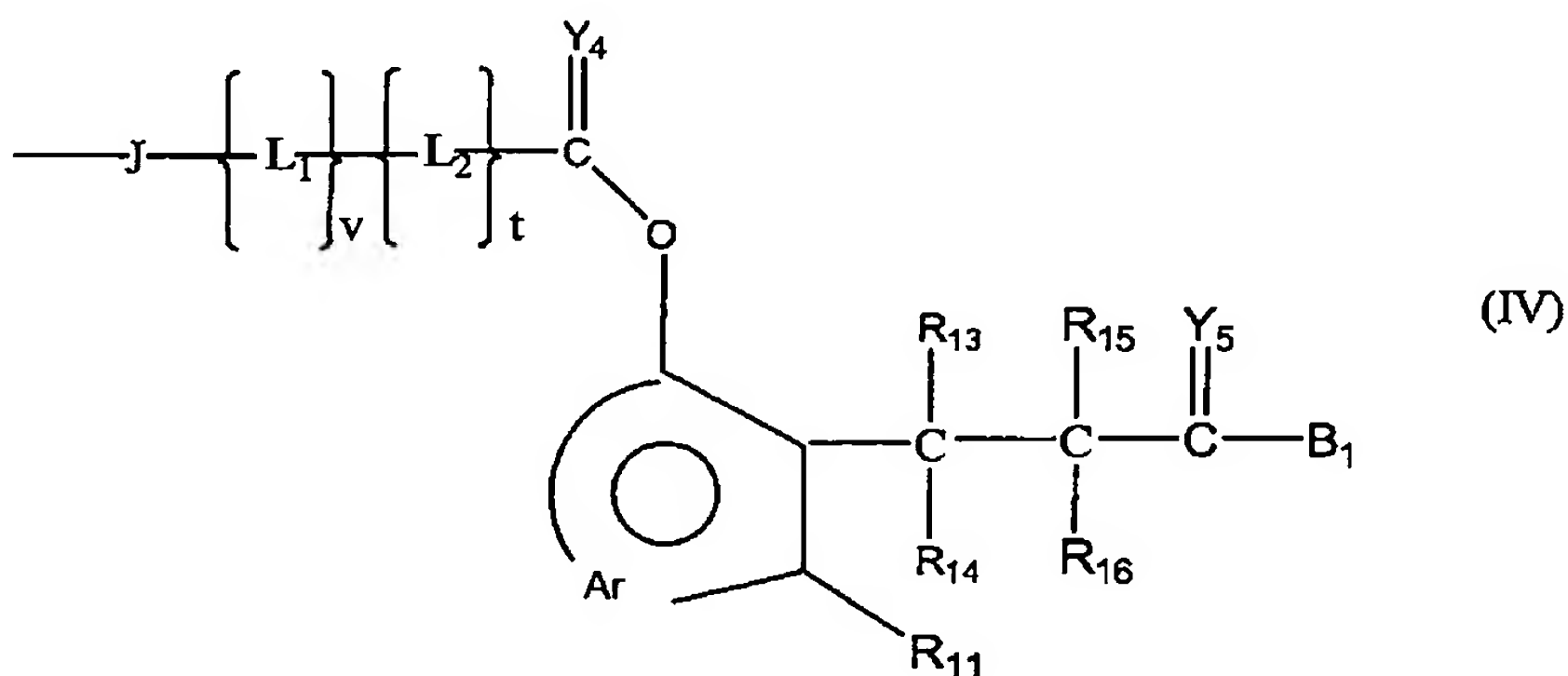


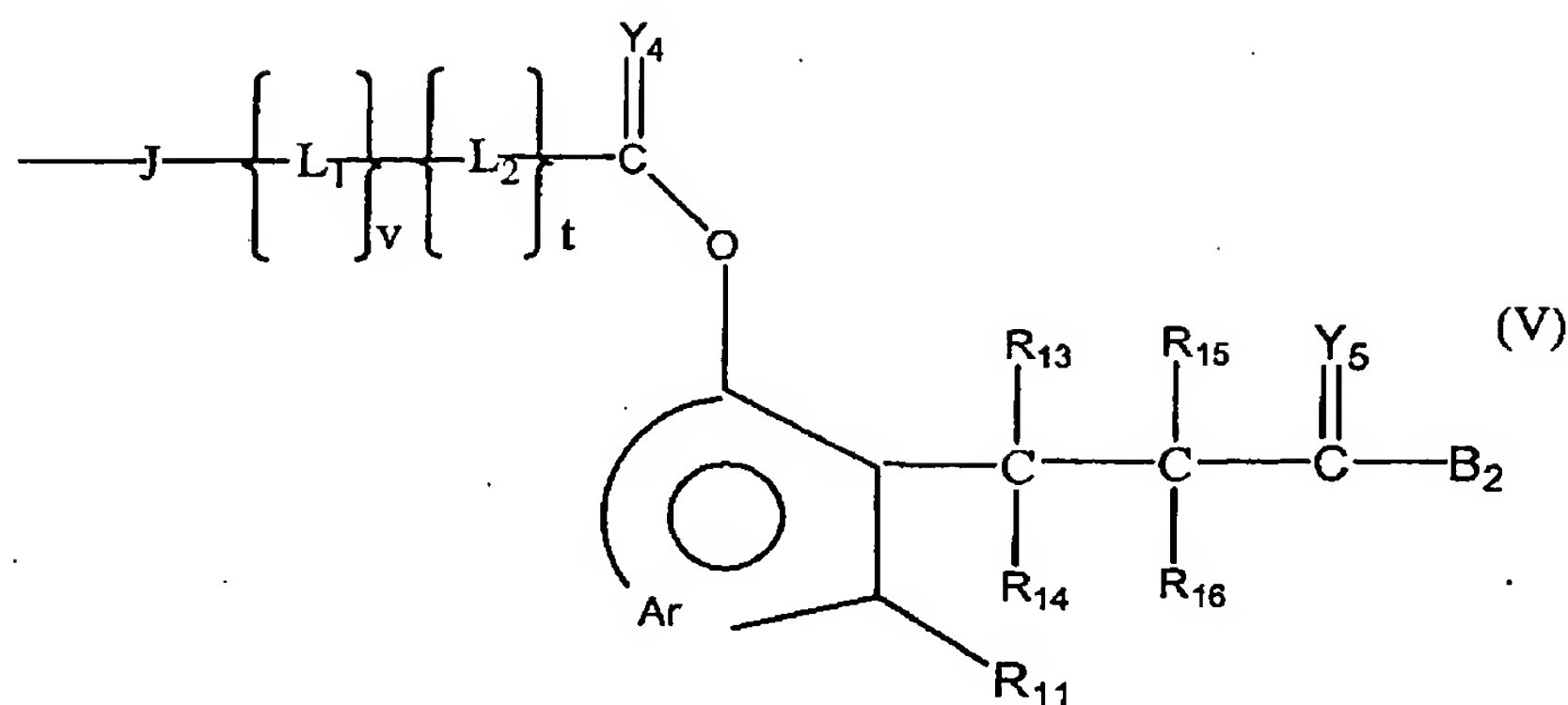
E_{35} is


$$\begin{array}{c} \text{R}_9 \\ | \\ \text{---C---} \\ | \\ \text{R}_8 \end{array} \begin{array}{c} \text{Y}_3 \\ || \\ \text{C} \end{array} \text{---D}'_2$$

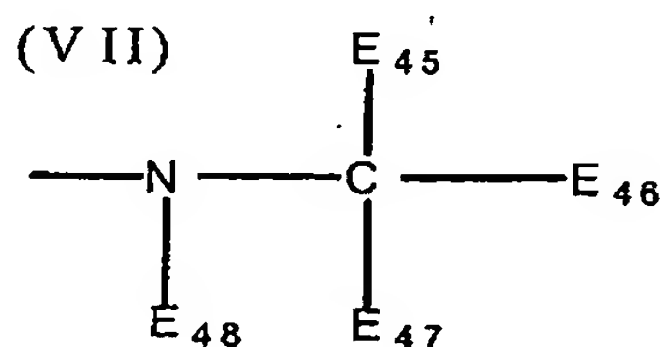
R_{6-10} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} hetero-alkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

D'_1 and D'_2 are independently OH ,





or



wherein (v) and (t) are independently 0 or a positive integer up to about 6;

L₁ and L₂ are independently selected bifunctional linkers;

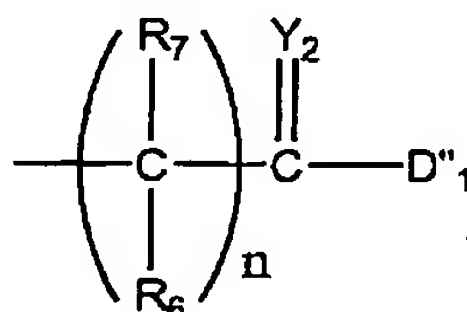
Y₄₋₇ are independently selected from the group consisting of O, S and NR₁₄;

R_{11-14} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} hetero-alkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

Ar is a moiety which when included in Formula (I) forms a multi-substituted aromatic hydrocarbon or a multi-substituted heterocyclic group;

B_1 and B_2 are independently selected from the group consisting of leaving groups, OH, residues of hydroxyl-containing moieties or amine-containing moieties;

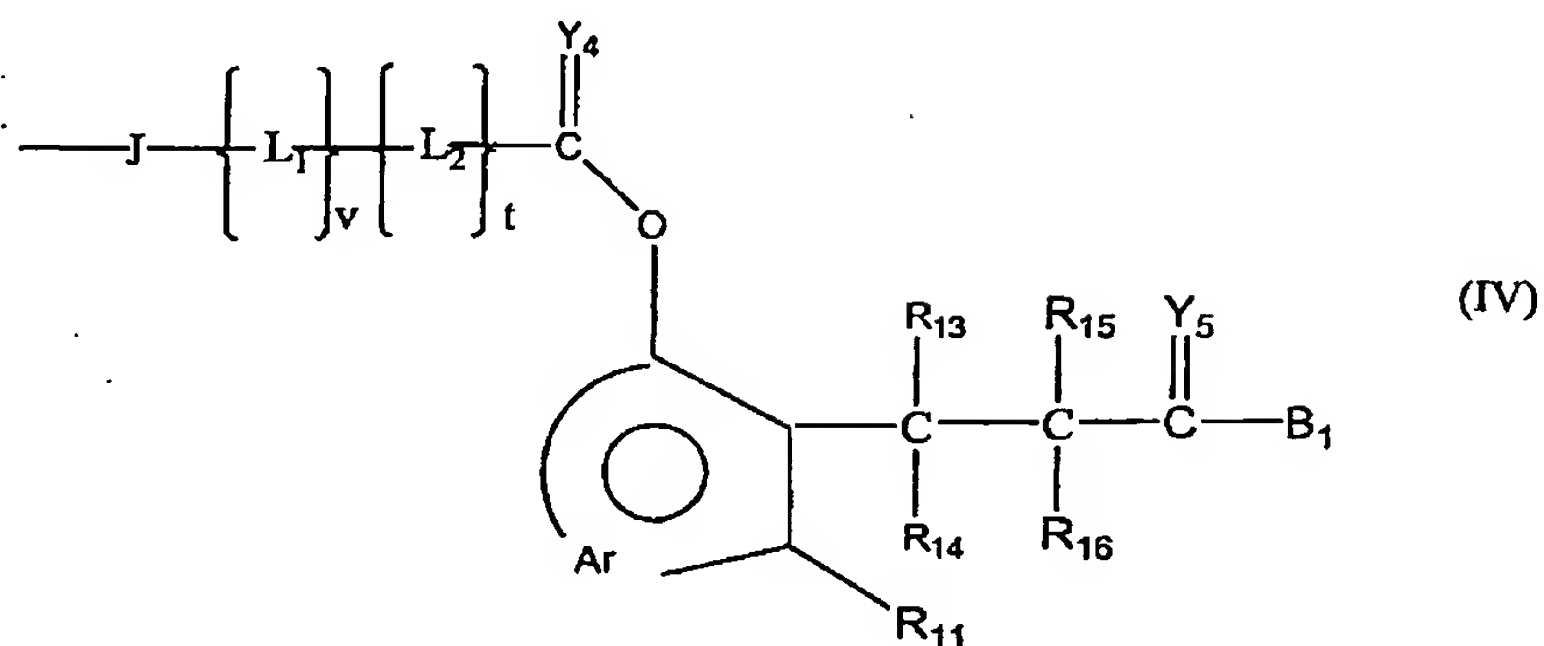
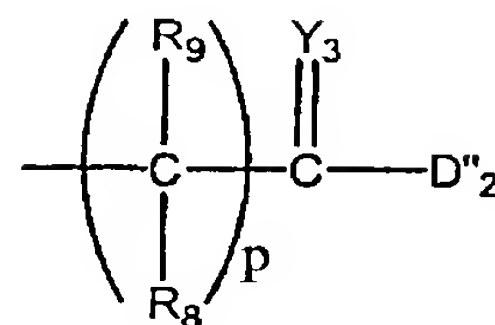
E_{45} is



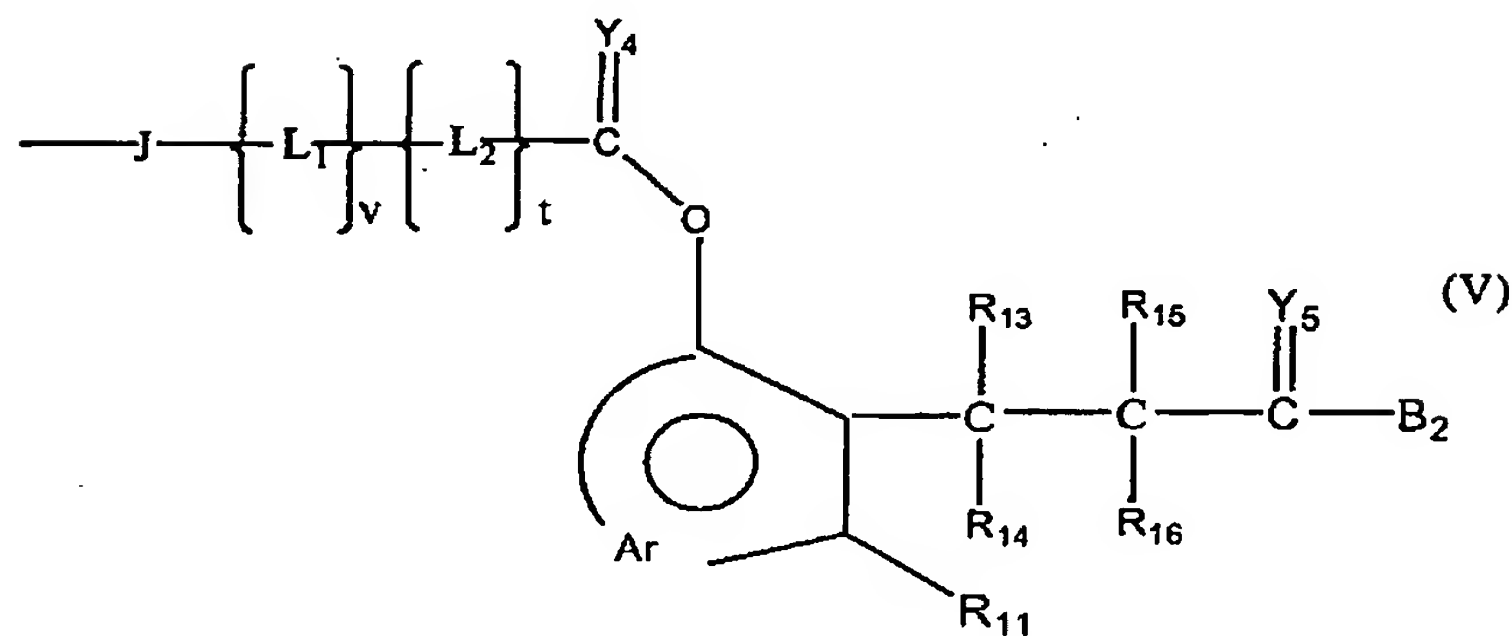
E_{46-48} are independently H, E_{45} or

wherein

D''_1 and D''_2 are independently OH,



or



5. (Currently amended) The compound of claim 3, wherein Y_1 is O.

6. (Original) The compound of claim 1, wherein R_1 comprises a polyalkylene oxide residue.

7. (Original) The compound of claim 6, wherein R_1 comprises a polyethylene glycol residue.

8. (Original) The compound of claim 3, wherein R_1 comprises a polyethylene glycol residue.

9. (Original) The compound of claim 6, wherein R_1 is selected from the group consisting of

$-C(=Y_6)-(CH_2)_f-O-(CH_2CH_2O)_x-A$,
 $-C(=Y_6)-Y_7-(CH_2)_f-O-(CH_2CH_2O)_x-A$,
 $-C(=Y_6)-NR_{23}-(CH_2)_f-O-(CH_2CH_2O)_x-A$,
 $-(CR_{24}R_{25})_e-O-(CH_2)_f-O-(CH_2CH_2O)_x-A$,
 $-NR_{23}-(CH_2)_f-O-(CH_2CH_2O)_x-A$,
 $-C(=Y_6)-(CH_2)_f-O-(CH_2CH_2O)_x-(CH_2)_f-C(=Y_6)-$,
 $-C(=Y_6)-Y_7-(CH_2)_f-O-(CH_2CH_2O)_x-(CH_2)_f-Y_7-C(=Y_6)-$,
 $-C(=Y_6)-NR_{23}-(CH_2)_f-O-(CH_2CH_2O)_x-(CH_2)_f-NR_{23}-C(=Y_6)-$,
 $-(CR_{24}R_{25})_e-O-(CH_2)_f-O-(CH_2CH_2O)_x-(CH_2)_f-O-(CR_{24}R_{25})_e-$, and
 $-NR_{23}-(CH_2)_f-O-(CH_2CH_2O)_x-(CH_2)_f-NR_{23}-$

wherein: Y_6 and Y_7 are independently O, S or NR_{23} ;

x is the degree of polymerization;

R_{23} , R_{24} and R_{25} are independently selected from among H, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

e and f are independently zero, one or two; and

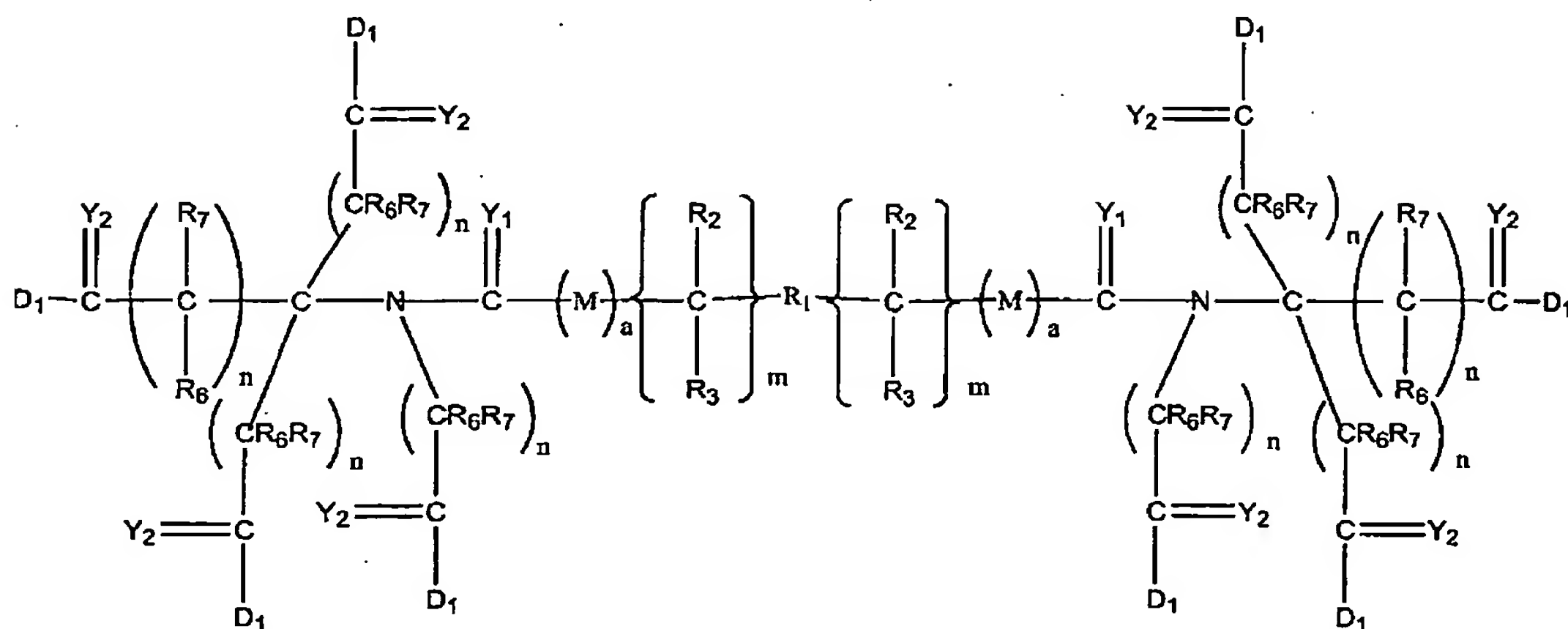
A is a capping group.

10. (Original) The compound of claim 9, wherein R_1 comprises $-O-(CH_2CH_2O)_x$ and x is a positive integer so that the weight average molecular weight is at least about 20,000.

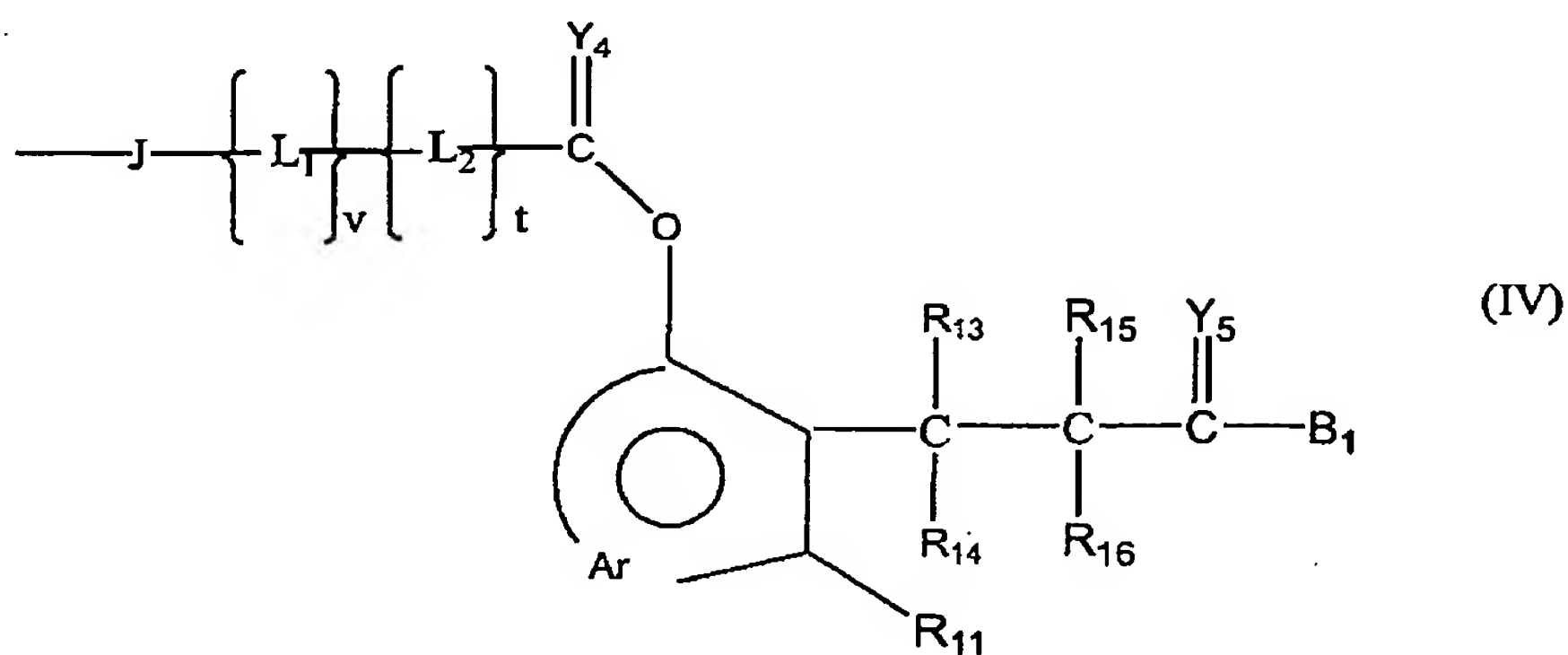
11. (Original) The compound of claim 3, wherein R_1 has a weight average molecular weight of from about 20,000 to about 100,000.

12. (Original) The compound of claim 3, wherein R₁ has a weight average molecular weight of from about 25,000 to about 60,000.

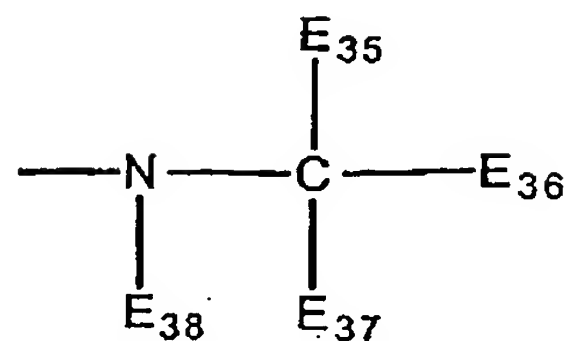
13. (Original) A compound of claim 3, comprising the formula



14. (Original) The compound of claim 13, wherein D₁ is



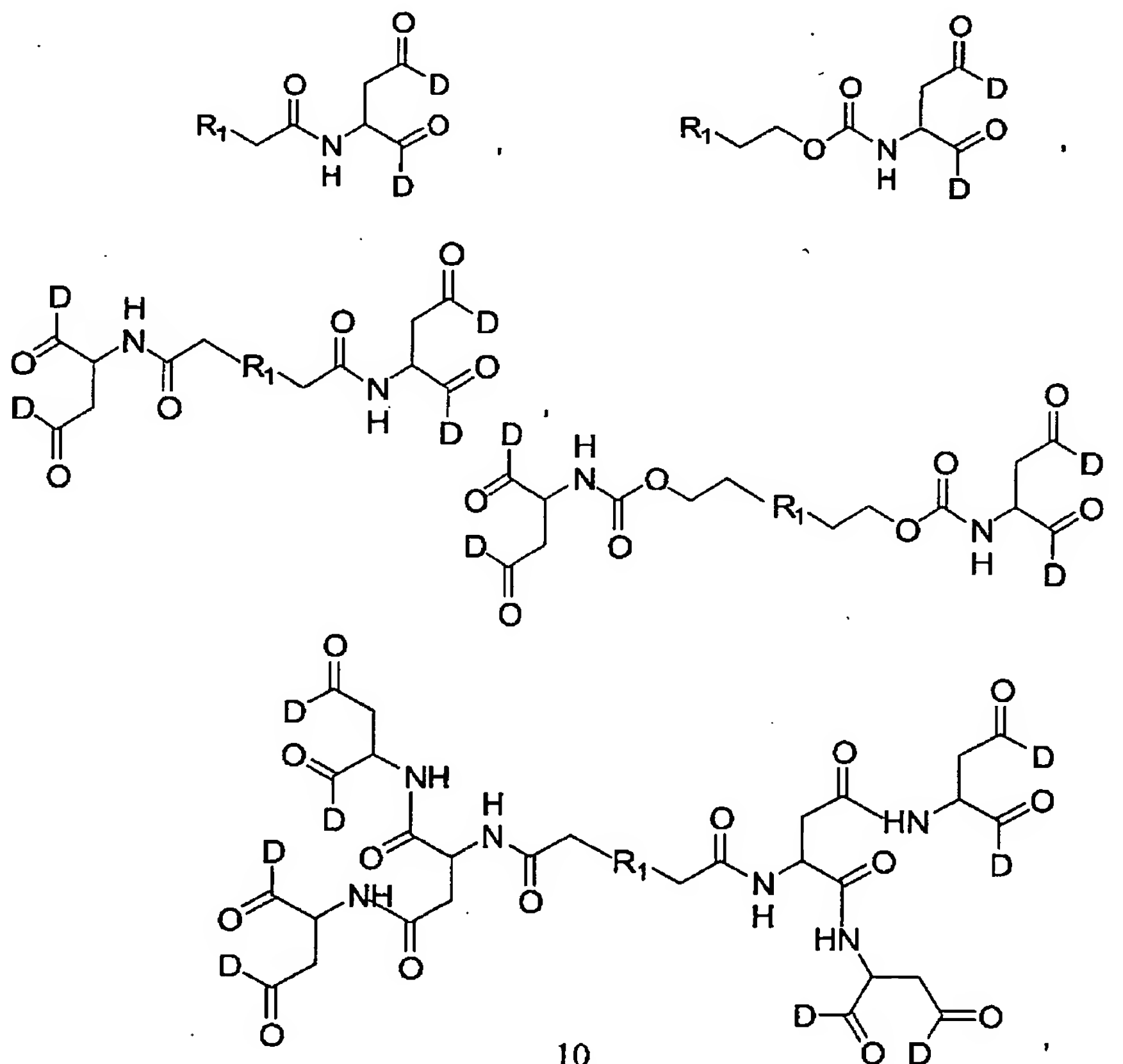
15. (Original) The compound of claim 13, wherein D_1 is

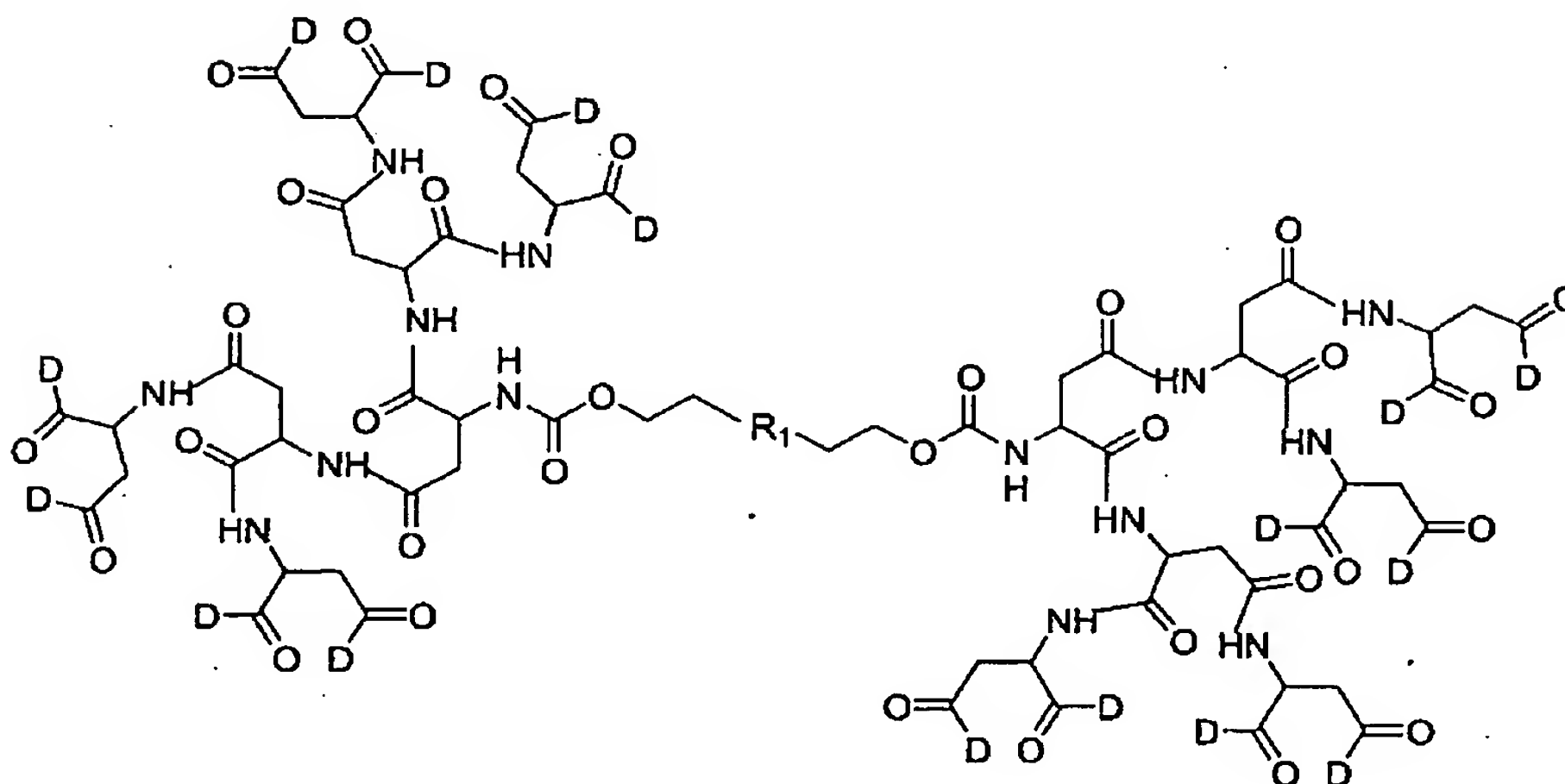
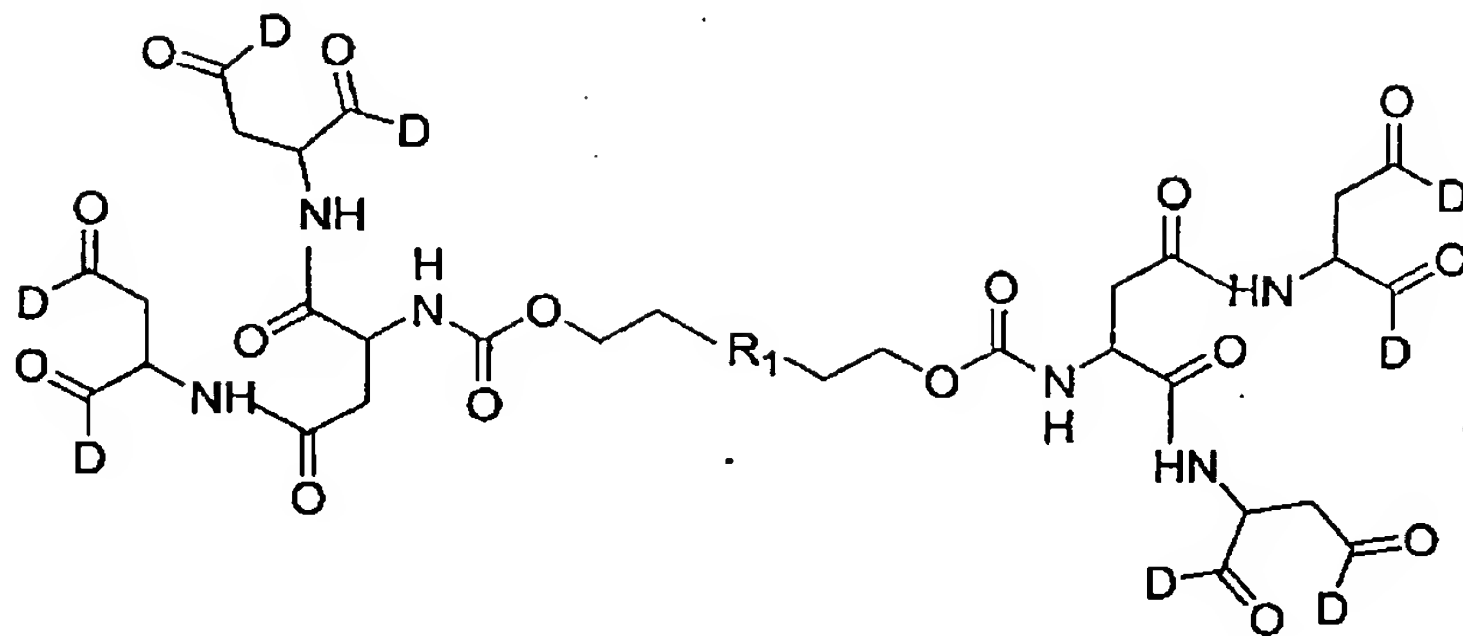


16. (Original) The compound of claim 1, wherein L_1 is $(CH_2CH_2O)_2$.

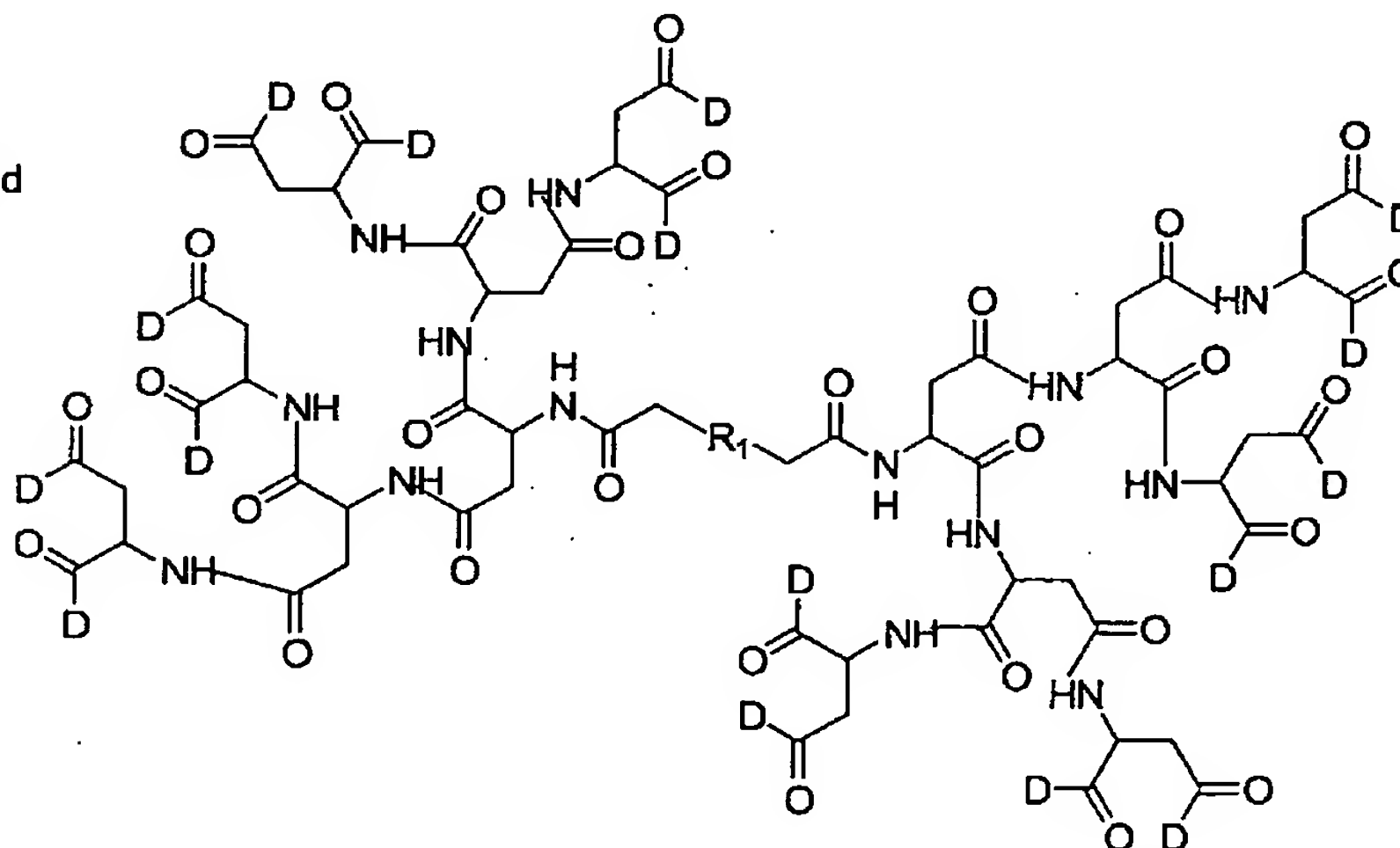
17. (Original) The compound of claim 1, wherein L_2 is selected from the group consisting of $-CH_2-$, $-CH(CH_3)-$, $-CH_2C(O)NHCH(CH_3)-$, $-(CH_2)_2-$, $-CH_2C(O)NHCH_2-$, $-(CH_2)_2-NH-$, $-(CH_2)_2-NH-C(O)(CH_2)_2NH-$ and $-CH_2C(O)NHCH(CH_2CH(CH_3)_2)-$.

18. (Original) A compound of claim 1, selected from the group consisting of:

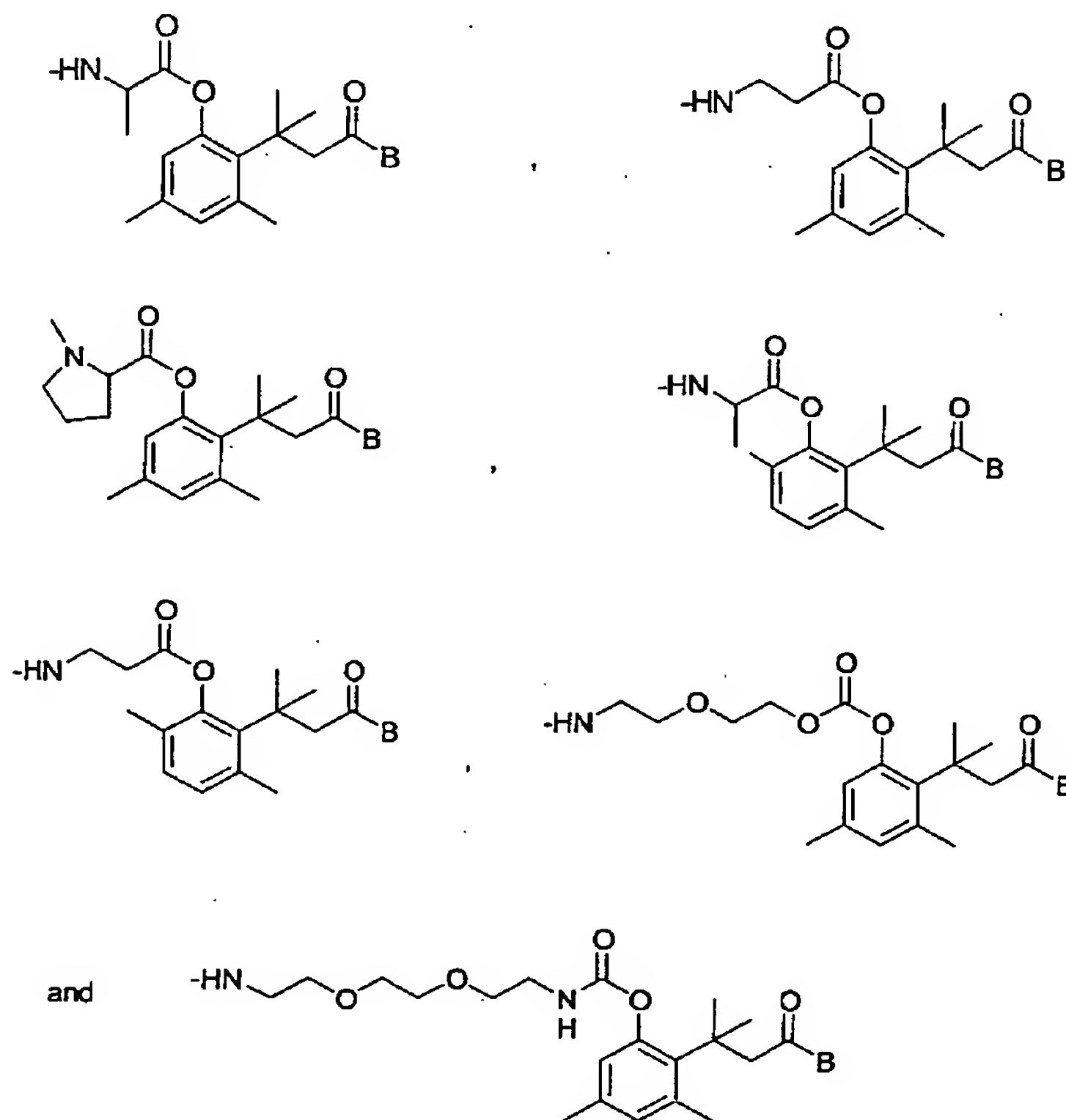




and



wherein R_1 is a PEG residue and D is selected from the group consisting of:



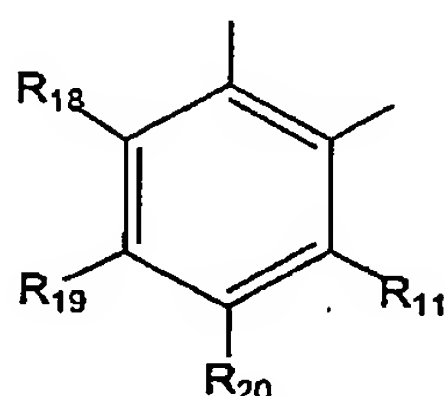
where B is a residue of an amine or a hydroxyl-containing drug.

19. (Original) A compound of claim 18, wherein B is a residue of a member of the group consisting of: daunorubicin, doxorubicin; *p*-aminoaniline mustard, melphalan, Ara-C (cytosine arabinoside), leucine-Ara-C, and gemcitabine

20. (Original) A method of treatment, comprising administering to a mammal in need of such treatment an effective amount of a compound of claim 1, wherein D_1 is a residue of a biologically active moiety.

21. (Original) A method of treatment, comprising administering to a mammal in need of such treatment an effective amount of a compound of claim 18.

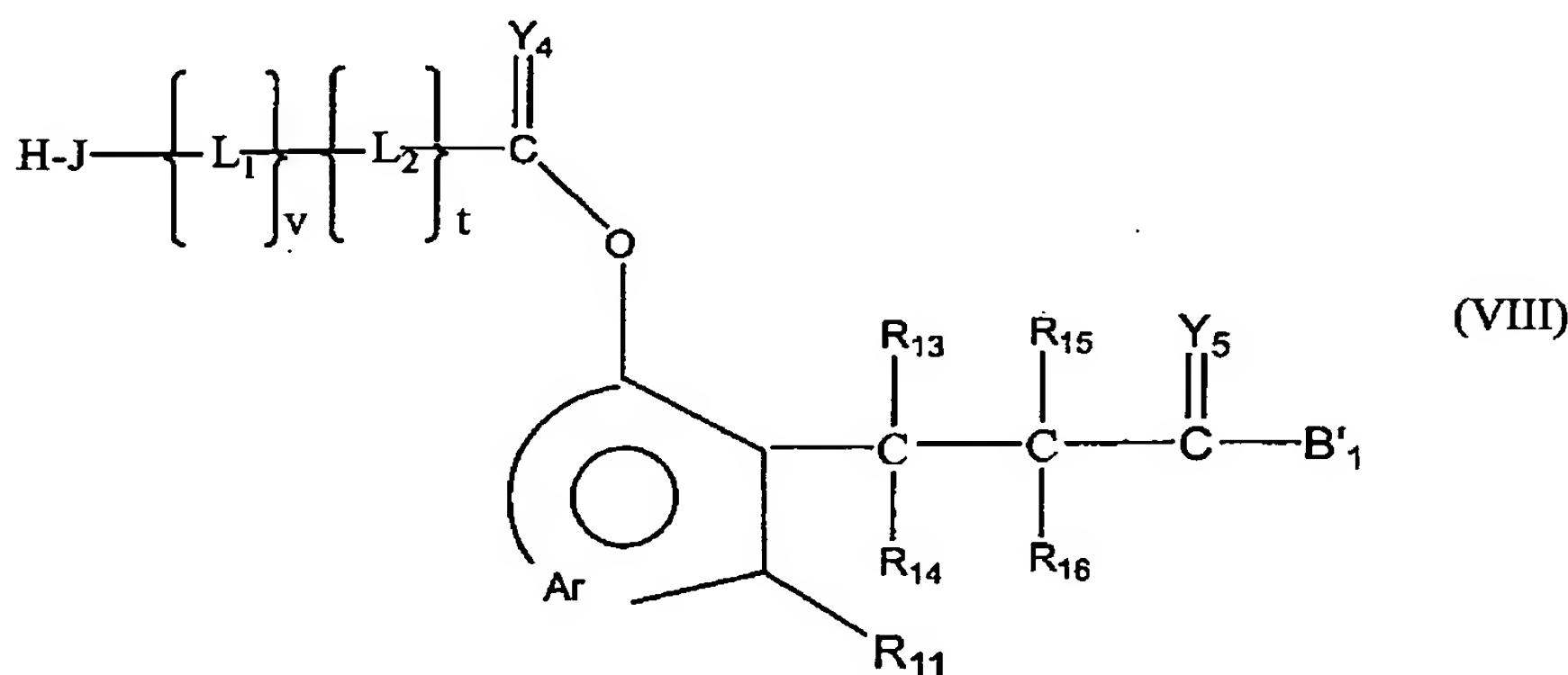
22. (Original) The compound of claim 1, wherein Ar comprises the formula:



wherein R₁₁ and R₁₈₋₂₀ are individually selected from the group consisting of hydrogen, C₁₋₆ alkyls, C₃₋₁₂ branched alkyls, C₃₋₈ cycloalkyls, C₁₋₆ substituted alkyls, C₃₋₈ substituted cycloalkyls, aryls, substituted aryls, aralkyls, C₁₋₆ heteroalkyls, substituted C₁₋₆ hetero-alkyls, C₁₋₆ alkoxy, phenoxy and C₁₋₆ heteroalkoxy.

23. (Original) The compound of claim 22, wherein R₁₁ and R₁₈₋₂₀ are each H or CH₃.

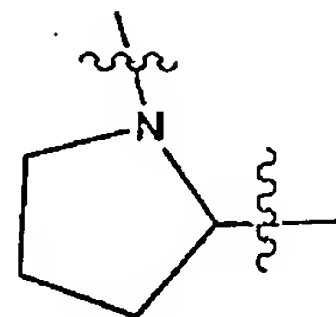
24. (Currently amended) A method of preparing a polymer conjugate, comprising:
reacting a compound of the formula (VIII):



wherein

(v) and (t) are independently 0 or a positive integer up to about 6;

J is NR_{12} or



L_1 and L_2 are independently selected bifunctional linkers;

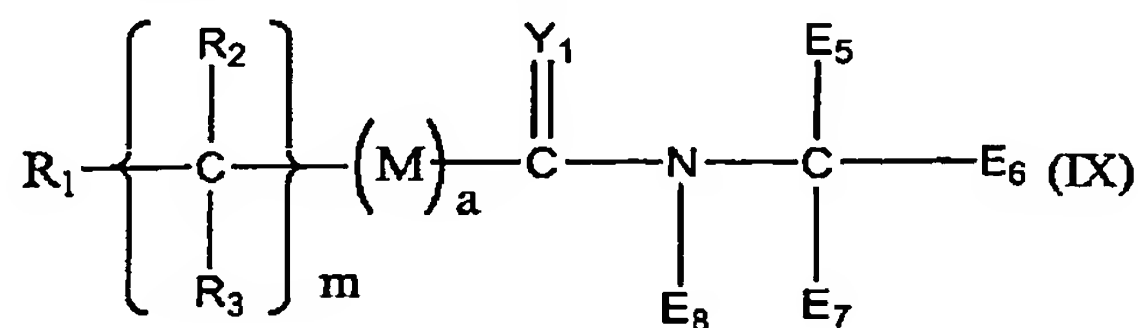
Y_{4-5} are independently selected from the group consisting of O, S and NR_{17} ;

R_{11-17} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} hetero-alkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

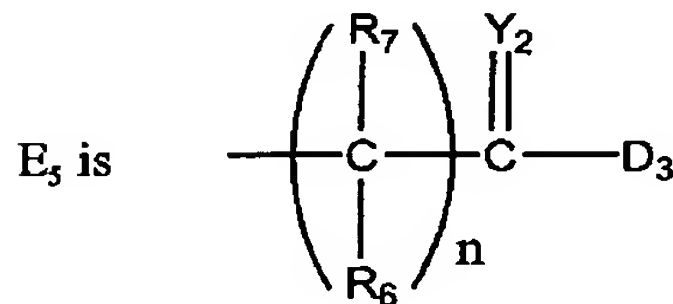
Ar is a moiety which when included in Formula (I) forms a multi-substituted aromatic hydrocarbon or a multi-substituted heterocyclic group; and

B' is a residue of a hydroxyl- or an amine-containing moiety;

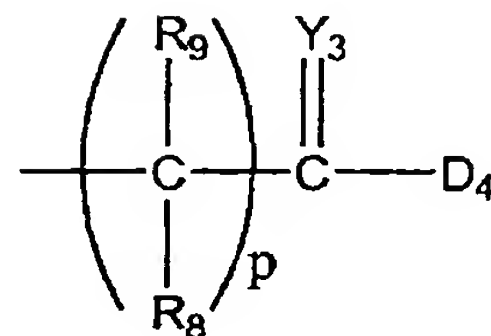
with a compound of the formula (IX):



wherein



E_{6-8} are independently H, E_5 or



D_3 and D_4 are independently OH, a leaving group which is capable of reacting with an unprotected amine or hydroxyl or a terminal branching group;

R_1 is a polymeric residue;

Y_1 is O, S or NR_4 ;

M is O, S or NR_5 ;

(a) is zero or one;

(m) is 0 or a positive integer;

(n) and (p) are independently 0 or a positive integer;

Y_{2-3} are independently O, S or NR_{10} ; and

R_{2-10} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} hetero-alkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

provided that E_{6-8} are not all H;

under conditions sufficient to cause a polymeric conjugate to be formed.